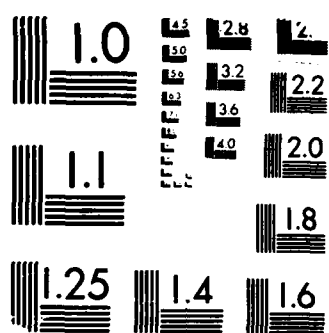


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A MINIMUM SPANNING PSEUDOFORREST

Harold N. Gabow  
Robert E. Tarjan

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July 1987

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# A LINEAR-TIME ALGORITHM FOR FINDING A MINIMUM SPANNING PSEUDOFORREST

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## **A Linear-Time Algorithm for Finding a Minimum Spanning Pseudoforest**

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### **Abstract.**

A *pseudoforest* is a graph each of whose connected components is a tree or a tree plus an edge; a *spanning pseudoforest* of a graph contains the greatest number of edges possible. This paper shows that a minimum cost spanning pseudoforest of a graph with  $n$  vertices and  $m$  edges can be found in  $O(m + n)$  time. This implies that a minimum spanning tree can be found in  $O(m)$  time for graphs with girth at least  $\log^{(i)} n$  for some constant  $i$ .

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## 1. Introduction.

A *pseudotree* is a connected graph with equal number of vertices and edges, i.e., a tree plus an edge creating a cycle. A *pseudoforest* is a graph each of whose connected components has at least as many vertices as edges, i.e., each component is a tree or a pseudotree. Pseudoforests arise in many applications although the terminology is not standard. We use the terminology of [PQ], which uses pseudoforests to compute the density and arboricity of a graph; see [W] for refinements of this approach. Pseudotrees are essentially the *1-trees* used in [HK] to solve the traveling salesman problem. The directed version of a pseudoforest is called a *functional graph* in [Be], since it corresponds to the graph of a finite function. For this reason pseudoforests commonly arise in parallel processing, when each processor chooses a successor (e.g., [GPS]). The pseudoforests of a graph form the *bicircular matroid*, which is important in the study of rigidity of bar-and-body frameworks [WW]. In the problem of minimum cost network flow with losses and gains [L], a linear programming basis is a pseudoforest [D]. A pseudotree is also called a *unicyclic graph* [e.g., MH].

With these applications as motivation we propose the *minimum spanning pseudoforest problem*: Consider a graph  $G$  with  $n$  vertices and  $m$  edges. A pseudoforest *spans*  $G$  if it has the greatest possible number of edges. Assume every edge  $e$  has a real-valued *cost*  $c(e)$ . The cost of a set of edges is the sum of all its edge costs. A *minimum spanning pseudoforest* has the smallest cost possible. This paper presents an algorithm to find such a pseudoforest in time  $O(m + n)$ .

The pseudoforest problem relates to finding a minimum spanning tree. The best-known time for finding a minimum spanning tree is  $O(m \log \beta(m, n))$  [GGST], where

$$\beta(m, n) = \min\{i \mid \log^{(i)} n \leq m/n\}.$$

Here  $\log$  denotes logarithm base two, and  $\log^{(i)} n$  is the  $i^{\text{th}}$  iterated logarithm, defined by  $\log^{(0)} n = n$ ,  $\log^{(i+1)} n = \log(\log^{(i)} n)$ . Note that if  $m/n \geq \log^{(i)} n$  for some constant  $i$  then  $\beta(m, n) \leq i$ , so the time to find a minimum spanning tree is  $O(m)$ . This paper presents a related result: If a graph has girth at least  $\log^{(i)} n$  for some constant  $i$  then a minimum spanning tree can be found in  $O(m)$  time.

Section 2 presents the results. This section closes with definitions and background from graph theory and data structures.

If  $S$  is a set and  $e$  an element,  $S + e$  denotes  $S \cup \{e\}$  and  $S - e$  denotes  $S - \{e\}$ . For a graph  $G$ ,  $V(G)$  and  $E(G)$  denote its vertex set and edge set, respectively. Hence for the given graph  $G$ ,  $n = |V(G)|$  and  $m = |E(G)|$ . An edge  $e$  is *incident* to a subgraph  $H$  if one or both ends is in  $V(H)$  but  $e \notin E(H)$ .

A *tree (pseudotree) component* of a graph  $G$  is a connected component of  $G$  that is a tree (pseudotree). A spanning pseudoforest  $P$  for a graph  $G$  consists of every tree component of  $G$ , plus for every other connected component  $C$  of  $G$ , one or more pseudotree components that partition  $V(C)$ . Note that  $P$  contains exactly  $|V(C)|$  edges of  $C$ .

The *set merging problem* [T] is to maintain a collection of disjoint sets which, after initialization, is subject to two operations:

*unite*( $S, S'$ )— form a new set  $S \cup S'$ , thereby destroying sets  $S$  and  $S'$ ;

*find*( $e$ )— return the name of the set containing element  $e$ .

The set merging algorithm used in Section 2 is *union by size*. It represents each set  $S$  by a *union tree*, i.e., a tree whose nodes are the elements of  $S$ . A *unite* makes the root of the smaller union tree a child of the root of the larger. An operation *find*( $v$ ) is done by following the path in the union tree from  $v$  to the root. (No path compression is done). Hence a *unite* operation is  $O(1)$  time and *find*( $v$ ) is  $O(\log s)$ , where  $s$  is the size of the set containing  $v$ .

In this paper a *priority queue* is a data structure on a universe that is partitioned into disjoint *queues*, where each element has a real-valued *cost*, and after initialization the following operations can be performed:

*meld*( $Q, Q'$ )— form a new queue by combining  $Q$  and  $Q'$ , thereby destroying queues  $Q$  and  $Q'$ ;

*find\_min*( $Q$ )— return the smallest cost element in queue  $Q$ ;

*delete*( $e, Q$ )— remove element  $e$  from queue  $Q$ .

The algorithm used in Section 2 implements priority queues with Fibonacci heaps [FT]. The following time bounds hold: *meld* is  $O(1)$ ; *find\_min*( $Q$ ) is  $O(\log s)$ , where  $s$  is the size of  $Q$ ; *delete*( $e, Q$ ) is  $O(\log s)$ , where  $s$  is the size of the Fibonacci tree containing  $e$ . Note these are amortized time bounds. Also to achieve the bound for *delete* the algorithm of [FT] is modified slightly, making it *lazier*: Unlike [FT] a queue does not keep track of its minimum element. Rather *find\_min*( $Q$ ) links trees of  $Q$  until there is at most one tree of each rank, and then finds and returns the desired minimum. *delete*( $e, Q$ ) cuts  $e$  from its parent and adds the children of  $e$  to the list of trees of  $Q$ . The analysis of [FT] easily extends to prove the above time bounds. (The same time bounds can be achieved using binomial queues [Br] modified to do lazy melding).

## 2. The algorithm.

The algorithm is based on a locality property similar to one possessed by minimum spanning trees [T].

**Lemma 2.1.** Let  $P$  be a subgraph of a minimum spanning pseudoforest. Let  $e$  be a smallest cost edge incident to some tree component  $T$  of  $P$ . Then  $P + e$  is a subgraph of a minimum spanning pseudoforest.

**Proof.** Let  $P^*$  be a minimum spanning pseudoforest containing  $P$ , and suppose  $P^*$  does not contain  $e$ . Let  $f$  be an edge of  $P^*$  that is incident to  $T$  such that the component of  $P^* - f$  containing  $T$  is a tree (Specifically if  $T$  is in a tree component of  $P$  then  $f$  is an edge of  $P$  incident to  $T$ ; if  $T$  is in a pseudotree component with cycle  $C$ , then  $f$  is an edge of  $P$  incident to  $T$  on  $C$  or on the path from  $T$  to  $C$ ). By definition,  $c(e) \leq c(f)$ . Hence  $P^* - f + e$  is the desired minimum spanning pseudoforest. ■

The algorithm enlarges a subgraph  $P$  to a minimum spanning pseudoforest. For efficiency it grows the components of  $P$  at approximately the same rate. More precisely let  $d(v)$  denote the degree of vertex  $v$  in the given graph  $G$ ; the (total) degree of a subgraph  $H$  is  $\sum\{d(v) | v \in V(H)\}$ . The algorithm grows components so that they have similar degrees. The details are as follows.

The algorithm initializes  $P$  to contain every vertex  $v$  of  $G$  ( $v$  is initially a tree component of  $P$ ). It then repeats the following step as long as  $P$  contains a tree component with an incident edge:

*Enlarging Step.* Choose a tree component  $T$  of smallest degree and add to  $P$  a minimum cost edge incident to  $T$ .

Correctness of this algorithm follows from the lemma; clearly pseudoforest  $P$  spans  $G$  when the algorithm halts.

The enlarging step is implemented with the following data structures. A set merging data structure maintains the partition of  $V(G)$  induced by the components of  $P$ . Each component of  $P$  is marked as a tree or pseudotree. Each tree component  $T$  maintains its degree  $d(T)$ , and a priority queue of incident edges  $Q(T)$ , ordered by cost. An edge can be in two priority queues, in which case the two occurrences are linked by pointers. There is an array  $C[1..2m]$ , where  $C[d]$  points to a doubly-linked list of all tree components of degree  $d$  with an incident edge.

With this data structure the enlarging step works as follows: The outermost loop examines the entries in  $C$  in increasing order to find the next smallest tree component  $T$ .  $T$  is removed from its  $C$ -list. The smallest edge  $e$  in  $Q(T)$  is obtained using *find\_min*. The set merging data structure finds the two components containing the ends of  $e$ , say  $T$  and  $S$ . If  $S = T$  it is marked



as a pseudotree. If  $S \neq T$  then sets  $V(S)$  and  $V(T)$  are united; further if  $S$  is a tree it is deleted from its  $C$ -list,  $e$  is deleted from  $Q(S)$  and  $Q(T)$ , these queues are merged, the new tree component  $S \cup T$  gets degree  $\delta = d(s) + d(t)$  and is added to the list  $C[\delta]$  if its queue is nonempty. Finally in all cases,  $e$  is added to  $P$ .

To estimate the time, note that all initialization uses  $O(m+n)$  time. The time for all enlarging steps, excluding priority queue *find\_mins* and *deletes* and set merging *finds*, is  $O(m+n)$ . To estimate the time for *find\_mins*, *deletes* and *finds*, define the rank of a component  $C$  as

$$r(C) = \lfloor \log d(C) \rfloor.$$

A simple induction shows that when  $T$  is chosen in the enlarging step, the size of any Fibonacci tree is at most  $d(T)$  (recall that *find\_min* is the only operation that enlarges Fibonacci trees; initially every edge is in its own Fibonacci tree). A similar induction shows that when  $T$  is chosen the height of the union tree for any component  $C$  is at most  $\min\{r(C), 1 + r(T)\}$  (since  $T$ 's height is at most  $r(T)$ ). Thus the *find\_min*, *find* and two *deletes* for  $T$  take time  $O(\log d(T) + r(T)) = O(r(T))$ . Let  $\mathcal{T}(r)$  denote the set of all rank  $r$  tree components chosen as  $T$  in the enlarging step. Then the total *find\_min*, *delete* and *find* time is at most a constant times

$$\sum_{r=0}^{\infty} r |\mathcal{T}(r)|.$$

For any rank  $r$ , any edge is counted in the degree of at most two trees of  $\mathcal{T}(r)$  (since the enlarging step unites  $T$  into a pseudotree or increases the rank of the component containing  $T$ ). Hence  $\sum \{d(T) | T \in \mathcal{T}(r)\} \leq 2m$ . Any  $T \in \mathcal{T}(r)$  has  $d(T) \geq 2^r$ . Thus  $|\mathcal{T}(r)| \leq m/2^{r-1}$ . This implies the total time is at most a constant times  $\sum_{r=0}^{\infty} rm/2^{r-1} = O(m)$ .

**Theorem 2.1.** A minimum spanning pseudoforest can be found in time  $O(m+n)$ . ■

Now we turn to the minimum spanning tree problem. Let  $P$  be a minimum spanning pseudoforest. Form a set  $C$  by choosing a maximum cost edge from each cycle of  $P$ .

**Lemma 2.2.**  $P - C$  is a subgraph of a minimum spanning tree.

**Proof.** Let  $T$  be a minimum spanning tree with as many edges of  $P$  as possible. Suppose  $P - C$  is not a subgraph of  $T$ . Let  $Q$  be a component of  $(P - C) \cap T$  that is not a component of  $P - C$ ; choose  $Q$  so it is not incident to an edge of  $C$ . Let  $e$  be an edge of  $P$  incident to  $Q$  such that the component of  $P - e$  containing  $Q$  is a tree ( $e$  is found as in Lemma 2.1). Let  $f$  be an edge incident



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